

=&gt; fil reg

FILE 'REGISTRY' ENTERED AT 13:41:57 ON 16 SEP 2002  
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.  
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.  
 COPYRIGHT (C) 2002 American Chemical Society (ACS)

*Structures 4a  
and  
5a*

Property values tagged with IC are from the ZIC/VINITI data file  
 provided by InfoChem.

STRUCTURE FILE UPDATES: 15 SEP 2002 HIGHEST RN 451445-11-7  
 DICTIONARY FILE UPDATES: 15 SEP 2002 HIGHEST RN 451445-11-7

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when  
 conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP  
 PROPERTIES for more information. See STNote 27, Searching Properties  
 in the CAS Registry File, for complete details:  
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=&gt; d his

(FILE 'REGISTRY' ENTERED AT 13:37:03 ON 16 SEP 2002)  
 DEL HIS Y  
 ACT FOUNDERED/A

L1 STR  
 L2 ( 6184)SEA FILE=REGISTRY SSS FUL L1  
 L3 STR  
 L4 5 SEA FILE=REGISTRY SUB=L2 SSS FUL L3

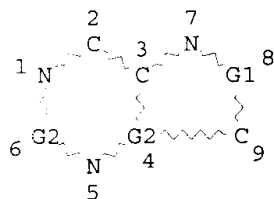
*→ none of these match  
claimed structures*

FILE 'REGISTRY' ENTERED AT 13:41:28 ON 16 SEP 2002

FILE 'REGISTRY' ENTERED AT 13:41:57 ON 16 SEP 2002

=&gt; d que stat l4

L1 STR



VAR G1=C/N/O/S

VAR G2=C/N

NODE ATTRIBUTES:

DEFAULT MLEVEL IS ATOM

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

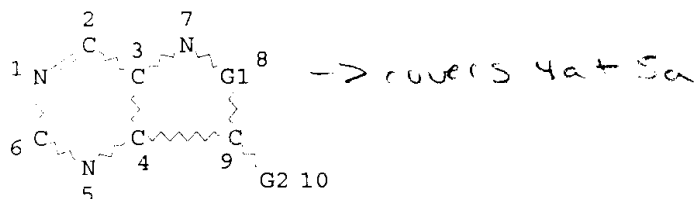
RSPEC I

NUMBER OF NODES IS 9

STEREO ATTRIBUTES: NONE

L2 ( 6184)SEA FILE=REGISTRY SSS FUL L1

L3 STR



VAR G1=O/S

VAR G2=AK/CY

NODE ATTRIBUTES:

DEFAULT MLEVEL IS ATOM

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 10

STEREO ATTRIBUTES: NONE

L4 5 SEA FILE=REGISTRY SUB=L2 SSS FUL L3

100.0% PROCESSED 89 ITERATIONS

5 ANSWERS

SEARCH TIME: 00.00.01

=&gt; d ide can l4 1-5

L4 ANSWER 1 OF 5 REGISTRY COPYRIGHT 2002 ACS

RN 173417-46-4 REGISTRY

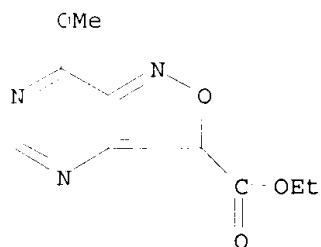
CN Isoxazolo[4,3-d]pyrimidine-3-carboxylic acid, 7-methoxy-, ethyl ester  
(9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C9 H9 N3 O4

SR CA

LC STN Files: CA, CAPLUS



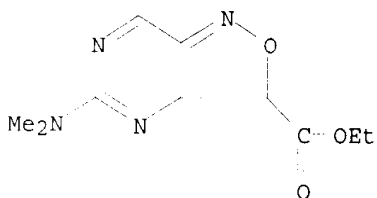
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 124:146101

L4 ANSWER 2 OF 5 REGISTRY COPYRIGHT 2002 ACS  
 RN 173417-45-3 REGISTRY  
 CN Isoxazolo[4,3-d]pyrimidine-3-carboxylic acid, 5-(dimethylamino)-, ethyl ester (9CI) (CA INDEX NAME)  
 FS 3D CONCORD  
 MF C10 H12 N4 O3  
 SR CA  
 LC STN Files: CA, CAPLUS, CASREACT

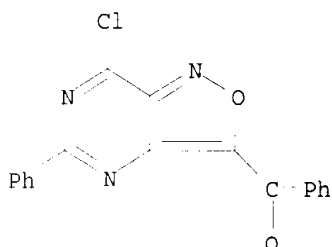


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1967 TO DATE)  
 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 124:146101

L4 ANSWER 3 OF 5 REGISTRY COPYRIGHT 2002 ACS  
 RN 166274-26-6 REGISTRY  
 CN Methanone, (7-chloro-5-phenylisoxazolo[4,3-d]pyrimidin-3-yl)phenyl- (9CI) (CA INDEX NAME)  
 FS 3D CONCORD  
 MF C18 H10 Cl N3 O2  
 SR CA  
 LC STN Files: CA, CAPLUS



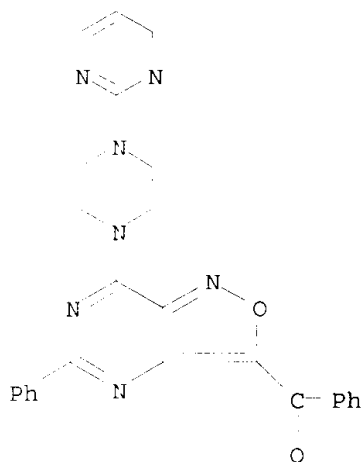
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1967 TO DATE)  
 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 123:143787

L4 ANSWER 4 OF 5 REGISTRY COPYRIGHT 2002 ACS  
 RN 166274-25-5 REGISTRY

CN Methanone, phenyl[5-phenyl-7-[4-(2-pyrimidinyl)-1-piperazinyl]isoxazolo[4,3-d]pyrimidin-3-yl]- (9CI) (CA INDEX NAME)  
 FS 3D CONCORD  
 MF C26 H21 N7 O2  
 SR CA  
 LC STN Files: CA, CAPLUS

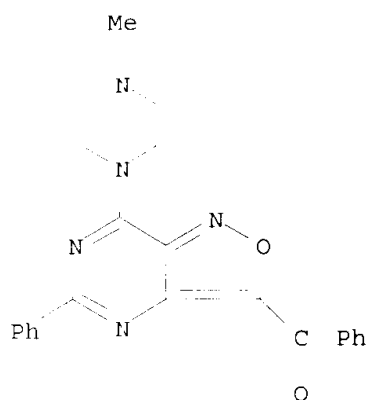


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1967 TO DATE)  
 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 123:143787

L4 ANSWER 5 OF 5 REGISTRY COPYRIGHT 2002 ACS  
 RN 166274-24-4 REGISTRY  
 CN Methanone, [7-(4-methyl-1-piperazinyl)-5-phenylisoxazolo[4,3-d]pyrimidin-3-yl]phenyl- (9CI) (CA INDEX NAME)  
 FS 3D CONCORD  
 MF C23 H21 N5 O2  
 SR CA  
 LC STN Files: CA, CAPLUS



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1967 TO DATE)  
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 123:143787